

In response to the Office Action of May 20, 2002, please amend the application as follows:

## IN THE SPECIFICATION

Between the title and "Field of Invention", insert the following

-- Cross Reference to Related Applications

This application is a national stage entry under 35 U.S.C. §371 of the PCT/EP99/05459 filed July 30, 1994.--

Page 1, lines 22-27:

Should rich as a should record

-(CH<sub>2</sub>),  $Ar_1$  where r is 0, 1 or 2 and Ar is an aromatic group chosen among benzene, naphthalene, thiophene, benzothiophene, pyridine, quinoline, indole, furan, benzofuran, thiazole, benzothiazole, imidazole, benzoimidazole, possibly substituted with up to 2 substituents chosen among  $C_{1-3}$  alkyl,  $C_{1-3}$  haloalkyl,  $C_{1-3}$  alkyloxy and  $C_{2-4}$  amino-alkyloxy, halogens, OH, NH<sub>2</sub>, CN, NR<sub>6</sub>R<sub>7</sub>, where R<sub>6</sub> and R<sub>7</sub>, are the same or different, and are H or  $C_{1-3}$  alkyl.

## Page 1, line 29-32 to Page 1a, line 2:

- $(CH_2)_rAr_1$  where r is 0, 1 or 2 and  $Ar_1$  is an aromatic group chosen among: benezene, naphthalene, thiophene, benzothiophene, pyridine, quinoline, indole, furan, benzofuran, thiazole, benzothiazole, imidazole, benzoimidazole, possibly substituted with up to 2 groups chosen among:  $C_{1-3}$  alkyl,  $C_{1-3}$ haloalkyl,  $C_{1-3}$  alkyloxy and  $C_{2-4}$  amino-alkyloxy, halogens, OH,  $NH_2$ , CN,  $NR_6R_7$ , where  $R_6$  and  $R_7$  are the same or different and are H or  $C_{1-3}$  alkyl.

Page 2, lines 10-12

or  $R_8$  and  $R_9$  together with the N atom to which they are linked to form a piperazine possibly substituted at one of its nitrogen atoms by  $C_{1-3}$  alkyl,  $C_{1-3}$  acyl or methanesulfonyl;

In WO9834949 it is described how compounds having lower molecular weight, monocyclic, containing only four bi-functional residues linked among each other by a peptide or pseudopeptide bond present pharmacological activity similar or higher than that of known compounds and moreover show a high selectivity for the human NK2 receptor.

Page 5, line 5 delete "triptophane" insert --tryptophane--.

More preferred are the compounds of formula (1) wherein:

- $-X_1$ ,  $X_2$ ,  $X_3$ ,  $X_4$  are -CONR-,
- -R is H;
- -R<sub>1</sub> is the lateral chain of tryptophane;

-R<sub>2</sub> is the lateral chain of phenylalanine substituted with up to two residues selected from the group consisting of: chlorine, fluorine, CF<sub>3</sub>, OH, and CN; or a group 3-pyridyl-methyl, or 4-pyridyl-methyl;

-R<sub>3</sub> is benzyl,

and the other substituents are as above defined.

## Page 5, lines 15-20

R<sub>9</sub> is a group chosen among: 4-tetrahydropyranyl, 4-tetraiodrothiopyranyl, 1-oxotetraiodrothiopyran-4-yl, 1,1 dioxo-tetrahydrothiopyran-4-yl, N-methyl-4-piperidinyl, N-methansulfonyl-4-piperidinyl, N-aminosulfonyl-4-piperdinyl, or R<sub>8</sub> and R<sub>9</sub> together with the N atom to which they are linked represent N-methyl-piperazinyl, N-acetyl-piperazinyl, piperazinyl, N-methanesulfonyl-piperazinyl.

## Page 10, line 32 and page 11, line 1

As starting compound the cyclo {Succinic[1-(R)-amino]-Trp-Phe-[(R)-NH-CH(CH $_2$ C $_6$ H $_5$ )-CH $_2$ -NH]-} (Compound A).

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